

The Low-Energy np Scattering Parameters and the Deuteron

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February 9, 2008

Abstract

The low-energy parameters describing the np scattering in the $^3S_1 + ^3D_1$ partial waves, the deuteron parameters, and their relations are discussed. These parameters can be determined quite accurately in the energy-dependent Nijmegen partial-wave analyses of the np data and are also given by the high-quality Nijmegen potentials. The newest values for these parameters are presented.

1 Introduction

The energy-dependent partial-wave analyses (PWA) of the NN scattering data as performed in Nijmegen [1] produce very good descriptions of these NN data below $T_L = 350$ MeV. We will consider explicitly the Nijm PWA93, which has $\chi^2/N_d = 0.99$ with respect to the scattering data [1]. The deuteron binding energy $B = 2.224575(9)$ MeV was included in these analyses as an experimental datum [2]. Because we have an energy-dependent solution of the scattering, we can produce accurate values for the low-energy scattering parameters and for the deuteron parameters. The PWA produces statistical errors on these parameters.

In Nijmegen we have also constructed high-quality (HQ) NN potentials [3]. These HQ potentials give a very good fit to the NN scattering data, including the deuteron binding energy. For each of these potentials we find in a direct comparison with the

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⁰Invited talk at the 3rd International Symposium “Dubna Deuteron 95”, Dubna, Moscow Region, Russia, July 4–7, 1995

NN scattering data that $\chi^2/N_d = 1.03$. Because the fit with these potentials is so good, we may speak here of alternative PWA's. Comparing the values of the low-energy parameters and the deuteron parameters as obtained in the various analyses gives us an estimate of the systematic error.

We can use these HQ potentials also to compute deuteron parameters that cannot be determined from the NN scattering data, such as the d -state probability p_d , the electric quadrupole moment Q , and the mean-square deuteron radius $\langle r^2 \rangle^{1/2}$. It turns out that also these parameters seem to be almost uniquely determined.

2 Scattering Matrix

The wave function is a solution of the relativistic Schrödinger equation

$$(\Delta + k^2)\psi = 2m V \psi ,$$

with m the reduced mass. The relativistic relation between the cm-momentum k and the cm-energy E is

$$E = \sqrt{k^2 + M_p^2} + \sqrt{k^2 + M_n^2} - (M_p + M_n) ,$$

where M_p and M_n are the proton and neutron mass. The asymptotic part of the wave function defines the partial-wave S matrix, which is related to the K matrix by

$$S = (1 + iK)/(1 - iK) .$$

For the spin-singlet waves (${}^1S_0, {}^1P_1, \dots$) and the uncoupled spin-triplet waves (${}^3P_0, {}^3P_1, {}^3D_2, \dots$), there is a simple relation between the phase shift δ and the S and K matrices, viz.

$$K = \tan \delta \quad \text{and} \quad S = e^{2i\delta} .$$

One can show that for low momenta $\delta_\ell \sim k^{2\ell+1}$.

For the triplet coupled waves (${}^3S_1 + {}^3D_1, {}^3P_2 + {}^3F_2, \dots$), there are two different parametrizations in use, the Stapp parametrization [4] and the Blatt and Biedenharn parametrization [5].

Stapp parametrization [4]

The (2×2) S matrix is written in terms of the nuclear-bar phase shifts $\bar{\delta}$ and $\bar{\varepsilon}$ as

$$S = \begin{pmatrix} e^{i\bar{\delta}_{J-1}} & \cdot \\ \cdot & e^{i\bar{\delta}_{J+1}} \end{pmatrix} \begin{pmatrix} \cos 2\bar{\varepsilon}_J & i \sin 2\bar{\varepsilon}_J \\ i \sin 2\bar{\varepsilon}_J & \cos 2\bar{\varepsilon}_J \end{pmatrix} \begin{pmatrix} e^{i\bar{\delta}_{J-1}} & \cdot \\ \cdot & e^{i\bar{\delta}_{J+1}} \end{pmatrix}$$

The advantage of this parametrization is that the behavior of the nuclear-bar mixing angle $\bar{\varepsilon}$ for low momenta is fine, in contradistinction with the behavior of the Blatt and Biedenharn [5] mixing angle ε (see below). For low momenta

$$\bar{\delta}_{J-1} \sim k^{2J-1}, \quad \bar{\delta}_{J+1} \sim k^{2J+3}, \quad \bar{\varepsilon}_J \sim k^{2J+1}.$$

Below the threshold $E = 0$ the analyticity properties of the nuclear-bar phases are not so nice.

Eigenphases (Blatt and Biedenharn) [5]

The (2×2) S matrix, and therefore also the K matrix, can be diagonalized by a real orthogonal matrix

$$O = \begin{pmatrix} \cos \varepsilon_J & -\sin \varepsilon_J \\ \sin \varepsilon_J & \cos \varepsilon_J \end{pmatrix},$$

with

$$S = O S_{\text{diag}} O^{-1} \quad \text{and} \quad K = O K_{\text{diag}} O^{-1}.$$

The eigenvalues define then the eigenphases δ_{J-1} and δ_{J+1} . The difficulty with this parametrization is that for low momenta

$$\delta_{J-1} \sim k^{2J-1}, \quad \delta_{J+1} \sim k^{2J+3}, \quad \text{but} \quad \varepsilon_J \sim k^2.$$

For $J > 1$ and small values of k the mixing parameter ε_J can have already a sizeable value, when the phases are both still practically zero. On the other hand, the analyticity properties of the eigenphases are nicer than of the nuclear-bar phases.

3 Analyticity

The S -matrix elements are analytic functions of the complex energy E or of the complex momenta k^2 . They have the so-called **unitarity cut** from $E = 0$ to ∞ along the positive real E axis. There are additional right-hand cuts starting at the pion-production threshold $E_{\text{cm}} \simeq 135$ MeV and left-hand cuts due to meson exchange. The OPE cut starts at $E = -m_\pi^2/(4M) = -5$ MeV, the TPE cut starts at $E = -m_\pi^2/M = -20$ MeV, where M = the nucleon mass and m_π = the pion mass. When there is a bound state present the S -matrix elements have a pole at the position of the bound state. In the ${}^3S_1 + {}^3D_1$ coupled channel there is therefore the deuteron pole at $E = -2.224575(9)$ MeV.

It is useful to define the **effective-range function** [6, 7]

$$M(k^2) = k^{\ell+1} K^{-1} k^\ell.$$

This effective-range function is again an analytic function of E or k^2 with right- and left-hand cuts as in the S -matrix elements. However, the unitarity cut is removed

and there is no pole at the position of the deuteron. The effective-range function $M(k^2)$ is therefore regular in a circle of radius 5 MeV around the origin, because then the OPE left-hand cut is reached. This effective-range function can be expanded in a power series around the origin with a radius of convergence of 5 MeV. One gets then the **effective-range expansion**

$$M = M_0 + M_1 k^2 + M_2 k^4 + M_3 k^6 + \dots$$

For practical purposes the radius of convergence of this effective-range expansion is pretty small (only 5 MeV). It is possible to define modified effective-range functions in which some of the left-hand cuts are removed, such as the OPE cut. The radius of convergence is then 20 MeV.

In the Nijmegen PWA we do not use the modified effective-range expansions [8, 9] anymore, but our methods, however, are quite similar.

4 The Deuteron in Scattering

The deuteron appears as a pole in the partial-wave S matrix for the coupled $^3S_1 + ^3D_1$ partial waves at $k = i\alpha$. The binding energy is given by

$$B = M_p + M_n - \sqrt{M_p^2 - \alpha^2} - \sqrt{M_n^2 - \alpha^2} = 2.224575(9) \text{ MeV} .$$

The radius R of the deuteron is then

$$R = 1/\alpha = 4.318946 \text{ fm} .$$

Here we use relativistic kinematics. The value of R in the case of non-relativistic kinematics is significantly different ($R_{\text{n.r.}} = 4.317667$ fm).

The partial-wave S matrix we write in the Blatt and Biedenharn parametrization

$$S = O \begin{pmatrix} S_0 & \cdot \\ \cdot & S_2 \end{pmatrix} O^{-1} .$$

The matrix elements of the matrix O are well behaved in the neighborhood of the deuteron pole. This allows us to define at the pole

$$\eta = -\tan \varepsilon_1 = 0.02543(7) .$$

The values quoted are from the Nijmegen 1993 PWA [1]. One of the eigenvalues (we take S_0) has a pole at $k = i\alpha$. We may write therefore

$$S_0 = \frac{N_p^2}{\alpha + ik} + (\text{regular function of } k) .$$

This expression is valid in the neighborhood of the deuteron pole. The **residue** at the pole is

$$N_p^2 = 0.7830(7) \text{ fm}^{-1} .$$

It is very convenient to define the effective range $\rho_d = \rho(-B, -B)$ at the deuteron pole by

$$\rho_d = R - 2/N_p^2 = 1.765(2) \text{ fm} .$$

Only 3 independent deuteron parameters B , η , and ρ_d can be determined from the np scattering data.

5 The Deuteron as a Bound State

The deuteron is a bound state in the coupled $^3S_1 + ^3D_1$ two-nucleon system. Let us neglect the possible existence of other coupled channels, such as $\Delta\Delta$, Q^6 , etc. In that case the deuteron wave function is

$$\psi = \frac{u(r)}{r} \mathcal{Y}_{011}{}^m + \frac{w(r)}{r} \mathcal{Y}_{211}{}^m ,$$

where $\mathcal{Y}_{LSJ}{}^m$ are the simultaneous eigenfunctions of the operators L^2 , S^2 , J^2 , and J_z . We assume the wave function to be properly normalized

$$\int_0^\infty dr (u^2 + w^2) = 1 .$$

The asymptotic behavior of the wave functions for large values of r are

$$u(r) \sim A_S e^{-\alpha r} \quad \text{and} \quad w(r) \sim A_D e^{-\alpha r} \left\{ 1 + 3(R/r) + 3(R/r)^2 \right\} .$$

The asymptotic normalizations are

$$A_S = 0.8841 \text{ fm}^{-1/2} \quad \text{and} \quad A_D = 0.0224 \text{ fm}^{-1/2} ,$$

with $N_d^2 = A_S^2 + A_D^2 = 0.7821 \text{ fm}^{-1}$. The d/s ratio $\eta = A_d/A_s = 0.02534$. The values quoted are from the Nijm I deuteron [3].

For the ordinary energy-independent local potentials and for velocity-dependent potentials one can show that $N_d = N_p$, i.e., the normalization as defined by the wave functions is equal to the normalization as defined by the residue [10]. In that case there is a trivial relation between the pole parameters determined in scattering and the asymptotic wave function. When we have energy-dependent potentials with $\partial U/\partial k^2 \neq 0$, then one can show that

$$(N_d/N_p)^2 = 1 - \int_0^\infty dr \left(\tilde{\psi}_d \frac{\partial U}{\partial k^2} \psi_d \right) .$$

6 Effective-Range Functions

The (2×2) matrix effective-range function is $M(k^2) = k^{\ell+1} K^{-1} k^\ell$. This matrix effective-range function can be expanded in a power series around the origin, but of course also around any other point in its analyticity domain, for example around the deuteron energy.

One can do the expansion around $k^2 = 0$, then

$$M(k^2) = M_0 + M_1 k^2 + M_2 k^4 + M_3 k^6 + M_4 k^8 + \dots ,$$

but one could do this also around $k^2 = -\alpha^2$, then

$$M(k^2) = m_0 + m_1(k^2 + \alpha^2) + m_2(k^2 + \alpha^2)^2 + \dots .$$

This means that

$$m_0 = M_0 - M_1 \alpha^2 + M_2 \alpha^4 - M_3 \alpha^6 + M_4 \alpha^8 - \dots$$

It is surprising, but also alarming, to discover that in order to get m_0 with sufficient accuracy one needs to include the term $M_6 \alpha^{12}$ in the right-hand side.

When we use the Blatt and Biedenharn parametrization in eigenphase shifts and coupling parameters, then we can define the effective-range functions

$$F_{J-1}(k^2) = k^{2J-1} \cot \delta_{J-1} , \quad F_{J+1}(k^2) = k^{2J+3} \cot \delta_{J+1} , \quad \text{and} \quad F_\varepsilon(k^2) = 2k^2 \cot 2\varepsilon_J .$$

These functions have the nice analyticity properties that make series expansions around the origin ($k^2 = 0$) or around the deuteron ($k^2 = -\alpha^2$) possible.

Using the Stapp parametrization we could define the effective-range functions

$$\bar{F}_{J-1} = 2k^{2J-1} \cot 2\bar{\delta}_{J-1} , \quad \bar{F}_{J+1} = 2k^{2J+3} \cot 2\bar{\delta}_{J+1} , \quad \text{and} \quad \bar{F}_\varepsilon = 2k^{2J+1} \cot 2\bar{\varepsilon}_J .$$

Again, the analyticity properties of these functions are not so nice.

7 Effective Ranges

It is customary to define various effective ranges such as the standard effective range $r = \rho(0, 0)$, the mixed effective range $\rho_m = \rho(0, -B)$, and the deuteron effective range $\rho_d = \rho(-B, -B)$. Let us look at the effective-range function $F_0(k^2) = k \cot \delta_0$ for the eigenphase δ_0 , also called the 3S_1 phase shift. Expanding in a power series around the origin gives

$$F_0(k^2) = -1/a + \frac{1}{2} r k^2 + v_2 k^4 + v_3 k^6 + \dots$$

where a = scattering length.

Some of the low-energy scattering parameters are given in Table 1. Remarkable is the agreement even in the parameters v_2 to v_5 . Numerical difficulties in our calculations prevented us to calculate accurately v_6 . We think that the values for v_5 are still

accurate, but we are not sure. That there is such an enormous agreement in these expansion coefficients comes in our opinion from our correct treatment of the OPE. It is this OPE that determines the fast energy dependence of the scattering matrix. The heavier-meson exchanges give rise to a much slower energy dependence.

It is easy to see that, because $S_0 = (1 + iK_0)/(1 - iK_0)$, one obtains a pole at $k^2 = -\alpha^2$ in the S -matrix elements when $K_0(-\alpha^2) = -i$. This implies that $F_0(-\alpha^2) = kK_0^{-1} = (i\alpha)(i) = -1/R$.

A series expansion around the deuteron is

$$F_0(k^2) = -1/R + \frac{1}{2}\rho_d(k^2 + \alpha^2) + w_2(k^2 + \alpha^2)^2 + \dots$$

The mixed effective range $\rho_m = \rho(0, -B)$ is defined by

$$-1/R = -1/a + \frac{1}{2}\rho_m(-\alpha^2) .$$

Then

$$\rho_m = 2R(1 - R/a) = 1.754(2)(3) \text{ fm} .$$

The first entry is the statistical error, the second one is the systematic error. We can make the expansions:

$$\begin{aligned} r &= \rho_m + 2\alpha^2 v_2 - 2\alpha^4 v_3 + 2\alpha^6 v_4 - \dots \\ \rho_d &= \rho_m - 2\alpha^2 v_2 + 4\alpha^4 v_3 - 6\alpha^6 v_4 - \dots \end{aligned}$$

The terms with v_6 in the above expansions are still significant. It turns out that

$$r \approx \rho_m - 0.001 \text{ fm} \quad \text{and} \quad \rho_d \approx \rho_m + 0.010 \text{ fm} .$$

Using the value for ρ_m gives

$$r = 1.753 \text{ fm} \quad \text{and} \quad \rho_d = 1.764 \text{ fm} .$$

At this point it becomes important to see how many independent low-energy parameters and deuteron parameters there are.

Table 1: Low-energy scattering parameters for the 3S_1 eigenphase.

	a	r	v_2	v_3	v_4	v_5
PWA	5.420(1)	1.753(2)	0.040	0.672	-3.96	27.1
Nijm I	5.418	1.751	0.046	0.675	-3.97	27.2
Nijm II	5.420	1.753	0.045	0.673	-3.95	27.0
Reid93	5.422	1.755	0.033	0.671	-3.90	26.7

Table 2: Deuteron parameters.

	η	ρ_d	p_d	$\langle r^2 \rangle^{1/2}$	Q_0
PWA93	0.02543	1.7647	–	–	–
Nijm I	0.02534	1.7619	5.664	1.9666	0.2719
Nijm II	0.02521	1.7642	5.635	1.9675	0.2707
Reid93	0.02514	1.7688	5.699	1.9686	0.2703

The binding energy B is directly related to the deuteron radius R . From the low-energy scattering follows the scattering length a . Using R and a gives the mixed effective range ρ_m . This determines then the values of the effective range r and the deuteron effective range ρ_d . This in turn determines the residue N_p^2 at the pole.

The d/s ratio η is also an independent quantity. We find therefore only 3 independent quantities: B , η , and a .

8 Multi-energy Solutions for S

As already pointed out in the introduction we have 4 excellent multi-energy parametrizations for the S matrix. The first one is our energy-dependent phase-shift analysis Nijm PWA93. We feel that the parameters determined in this analysis [1] are probably the most accurate. Then there exist 3 HQ potentials [3] (Nijm I, Nijm II, and Reid93) which can be considered as alternative PWA's with a slightly higher χ^2_{\min} than Nijm PWA93. In Table 1 we presented already some of the effective-range parameters. In Table 2 we present the various deuteron parameters. The deuteron parameters η and ρ_d can be determined in PWA's as well as from the HQ potentials. The d -state probability p_d , the mean-square radius $\langle r^2 \rangle^{1/2}$, and the electric quadrupole moment can only be determined for the HQ potentials.

9 Conclusions

In this contribution we presented the various low-energy parameters and the deuteron parameters. We would like to stress once more that effective-range expansions are not very accurate, unless one takes many terms into account. A similar situation exists in the 1S_0 channel. Also there the effective-range expansion is not accurate enough to describe the data with sufficient accuracy [9]. As a result, the effective-range parameters have only a very limited usefulness. The following low-energy parameters and deuteron parameters we would like to recommend. The errors are educated guesses for the total error, including statistical as well as systematic errors.

- scattering length $a = 5.4194(20)$ fm.
- effective range $r = 1.7536(25)$ fm.
The values of a and r are strongly correlated due to the accurate value of the deuteron binding energy. A large value of a implies a large value of r .
- shape parameters $v_2 = 0.040(7)$ fm³, $v_3 = 0.673(2)$ fm⁵, $v_4 = 3.95(5)$ fm⁷, and $v_5 = 27.0(3)$ fm⁹.
- deuteron binding energy $B = 2.224575(9)$ MeV.
- deuteron radius (rel) $R = 4.318946$ fm.
- d/s -ratio $\eta = 0.0253(2)$.
This value is in good agreement with the recent determination by Rodning and Knutson [11] of $\eta = 0.0256(4)$.
- deuteron effective range $\rho_d = 1.765(4)$ fm.
- residue $N_p^2 = 0.7830(15)$ fm.
- asymptotic normalizations $A_S = 0.8845(8)$ fm^{-1/2} and $A_D = 0.0223(2)$ fm^{-1/2}.
- d -state probability $p_d = 5.67(4)\%$.
- mean-square radius $\langle r^2 \rangle^{1/2} = 1.9676(10)$ fm.
This must be compared with the recent experimental determination in atomic physics [12], which results in: $\langle r^2 \rangle^{1/2} = 1.971(6)$ fm. For good discussions see Refs [13, 14].
- electric quadrupole moment $Q_0 = 0.271(1)$ fm².
This value disagrees with the experimental value [15] $Q = 0.2859(3)$ fm². The difference $Q - Q_0 \simeq 0.015$ fm² needs to be explained in terms of meson exchange currents, relativistic effects, contributions of Q^6 states, etc.

We would like to point out here the existence of the NN-OnLine facility of the Nijmegen group [16], which is accessible via the World-Wide Web. There one can obtain the various Nijmegen e-prints, the fortran codes for some of the Nijmegen potentials, the deuteron parameters and the deuteron wave functions, the phases obtained from the Nijmegen PWA and from the Nijmegen potentials, and predictions for many of the experimental quantities. A direct comparison of these predictions with the Nijmegen NN-data base is also possible.

Acknowledgments

We would like to thank the other members of the Nijmegen group for their interest, help, and discussions.

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